## **Discretization Methods**

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- Consider an AR(1) process of the form

$$z_t = \rho z_{t-1} + \varepsilon_t \tag{1}$$

with stationary distribution  $\mathcal{N}(0, \sigma_z^2)$  where  $\sigma_z = \frac{\sigma_{\varepsilon}}{\sqrt{1-\rho^2}}$ 

- **Goal**: obtain a finite state Markov chain that generates the same population moments as the continuous process
- Why useful?
  - \* Solving Bellman equations typically involves taking expectations over next period values
  - \* If process is continuous, then one has to deal with integrals which are computationally costly
  - \* If discrete, the continuation value is just a weighted sum



- Let *X* be a finite set with *n* elements :  $\{x_1, \ldots, x_n\}$ 
  - \* Think of *X* as the *exogenous* state space and *x<sub>i</sub>* as the state values
- A Markov chain on X is a sequence of random variables {X<sub>t</sub>} that satisfy for any date t and any next period state x̃ ∈ X

$$\mathbb{P}(X_{t+1} = \tilde{x} | X_t) = \mathbb{P}(X_{t+1} = \tilde{x} | X_t, X_{t-1}, X_{t-2}, \ldots)$$
(2)

- \* In words, knowing the current state is enough to form expectations about future states.
- \* This is the so called Markov property
- The dynamics of a Markov chain are fully determined by the set of probability values

$$P(x,\tilde{x}) := \mathbb{P}(X_{t+1} = \tilde{x} | X_t = x) \quad (x,\tilde{x} \in X)$$
(3)

- \*  $P(x, \tilde{x}) \ge 0$  is the probability of going from x to  $\tilde{x}$  in one step
- \*  $P(x, \cdot)$  is the conditional distribution of  $X_{t+1}$  given  $X_t = x$



- Assume that you have a stochastic matrix P, that is a  $n \times n$  matrix such that
  - \* Each element of *P* is non-negative
  - \* Each row of *P* sums to one
- Then, you can generate a Markov chain  $\{X_t\}$  as follows:
  - \* Set a initial value or draw it from a know distribution. Call it  $x_0$ .
  - \* Since you know  $x_0$ , you can draw  $X_1$  using the  $\mathbb{P}(X_1|X_0 = x_0)$
  - \* Thus, in general you can draw  $X_{t+1}$  using  $P(X_t, \cdot)$  for all t = 0, 1, ...
- Let's see an example using the following stochastic matrix

$$\boldsymbol{P} := \begin{bmatrix} \alpha & 1-\alpha \\ 1-\beta & \beta \end{bmatrix}$$
(4)

\* You can think of  $1 - \alpha$  as the probability of finding a job conditional on being unemployed and  $1 - \beta$  as the probability of losing a job conditional on being employed.



1. Choose a value for the persistence  $ho \in (0, 1)$  and the standard deviation  $\sigma_z > 0$ 

- 2. Set values for the hyper-parameters
  - \* *n* : number of potential realization of the process
  - \* *m* : number of standard deviations away from the the unconditional mean
  - \* Typical values : n= $\{5, 7, 9, 11, 15\}, m = \{2, 3, 4\}$
- 3. Set the bounds for the process

$$\bar{z} = m\sigma_z$$
(5)
 $z = -m\sigma_z$ 
(6)

## Tauchen's Method



4. Set  $\{z_i\}_{i=1}^n$  such that:

$$z_i = \underline{z} + \frac{\overline{z} - \underline{z}}{n-1}(i-1)$$
(7)

and construct mid-points  $\{\tilde{z}\}_{i=1}^{n-1}$  which are given by:

$$\tilde{z}_i = \frac{z_{i+1} + z_i}{2} \tag{8}$$

5. The transition probability  $p_{ij} \in P_{z,z'}$  (the probability of going to state  $z_j$  conditional on being in state  $z_i$ ) is computed according to

$$p_{ij} = \Phi\left(\frac{\tilde{z}_j - \rho z_i}{\sigma_{\varepsilon}}\right) - \Phi\left(\frac{\tilde{z}_{j-1} - \rho z_i}{\sigma_{\varepsilon}}\right) \quad j = 2, 3, \dots, n-1$$
(9)

$$p_{i1} = \Phi\left(\frac{\tilde{z}_1 - \rho z_i}{\sigma_{\varepsilon}}\right) \tag{10}$$

$$p_{in} = 1 - \Phi\left(\frac{\widetilde{z}_{n-1} - \rho z_i}{\sigma_{\varepsilon}}\right) \tag{11}$$

where  $\Phi(\cdot)$  denotes a CDF of the  $\mathcal{N}(\mathbf{0},\mathbf{1})$ 



- Expression (9) is obtained as follows
  - \* Let  $d = z_{k+1} z_k$  be the distance between two points in the vector of state values.

\* Then,

$$p_{i,j} = \Pr \{ z' = z_j \mid z = z_i \}$$

$$= \Pr \{ z_j - d/2 < z' \le z_j + d/2 \mid z = z_i \}$$

$$= \Pr \{ z_j - d/2 < \rho z_i + \varepsilon \le z_j + d/2 \}$$

$$= \Pr \{ \frac{z_j + d/2 - \rho z_i}{\sigma_{\varepsilon}} < \frac{\varepsilon}{\sigma_{\varepsilon}} \le \frac{z_j - d/2 - \rho z_i}{\sigma_{\varepsilon}} \}$$

$$= \Phi \left( \frac{z_j + d/2 - \rho z_i}{\sigma_{\varepsilon}} \right) - \Phi \left( \frac{z_j - d/2 - \rho z_i}{\sigma_{\varepsilon}} \right)$$
(12)



- Assume that we have a stochastic process

$$y_t = 0.85y_{t-1} + \varepsilon$$
 with  $\mathcal{N}(0, 0.0095^2)$  (13)

- We want to approximate it with a Markov chain with 5 and 9 points. We set m = 3.
- How does the approximation depend on number of potential realizations of the process?
  - \* We simulate the Markov chain for T = 10,000 periods
  - \* Are the sample moments close to the population ones?
  - \* What about the persistence of the process?

|       | $\hat{\mu}_{y}$ | $\hat{\sigma}_{y}$ | ρ     |  |
|-------|-----------------|--------------------|-------|--|
| n = 5 | 1.42e-5         | 0.012              | 0.881 |  |
| n = 9 | -0.0001         | 0.010              | 0.851 |  |



- 1. Choose a value for the persistence  $ho\in(0,1)$  and the standard deviation  $\sigma_z>0$
- 2. Set values for the hyper-parameters
  - \* *n* : number of potential realization of the process
  - \*  $\lambda$  : controls the upper and lower bound of the process
  - \* Typical value for  $\lambda = \sqrt{n-1} \sigma_z$
- 3. Set the bounds for the process as follows

$$\bar{z} = \lambda$$
 (14)

$$\underline{z} = -\lambda$$
 (15)

4. Set  $\{z_i\}_{i=1}^n$  such that:

$$z_i = \underline{z} + \frac{\overline{z} - \underline{z}}{n-1}(i-1)$$
(16)

## Rouwenhorst's Method



5. When n = 2, let  $P_2$  be given by

$$P_2 = \begin{bmatrix} p & 1-p \\ 1-q & q \end{bmatrix}$$

\* *p* and *q* can be set to  $p = q = \frac{1+\rho}{2}$ 

6. For n > 2, construct *recursively* the transition matrix:

$$P_{n} = p \begin{bmatrix} P_{n-1} & \mathbf{0} \\ \mathbf{0}' & \mathbf{0} \end{bmatrix} + (1-p) \begin{bmatrix} \mathbf{0} & P_{n-1} \\ \mathbf{0} & \mathbf{0}' \end{bmatrix} + q \begin{bmatrix} \mathbf{0}' & \mathbf{0} \\ P_{n-1} & \mathbf{0} \end{bmatrix} + (1-q) \begin{bmatrix} \mathbf{0} & \mathbf{0}' \\ \mathbf{0} & P_{n-1} \end{bmatrix}$$
(18)

where **0** is a  $(n-1) \times 1$  column vector of zeros.

7. Divide all elements in the middle rows (except top and bottom) by 2 so the sum of each row is equal to 1. The final outcome is  $P_{z,z'}$ 

(17)





- Let n = 3 and  $P_2$  be given by equation (17). Then,

$$P_{3} = p \begin{bmatrix} p & 1-p & 0\\ 1-q & q & 0\\ 0 & 0 & 0 \end{bmatrix} + (1-p) \begin{bmatrix} 0 & p & 1-p\\ 0 & 1-q & q\\ 0 & 0 & 0 \end{bmatrix} + q \begin{bmatrix} 0 & 0 & 0\\ p & 1-p & 0\\ 1-q & q & 0 \end{bmatrix} + (1-q) \begin{bmatrix} 0 & 0 & 0\\ 0 & p & 1-p\\ 0 & 1-q & q \end{bmatrix}$$

- After multiplying and re-arranging terms we obtain

$$P_{3} = \begin{bmatrix} p^{2} & 2p(1-p) & (1-p)^{2} \\ p & 1 & 1-p \\ (1-q)q & q^{2} + (1-q)^{2} & (1-q)q \end{bmatrix}$$
(20)

- Second row sums up to 2! Not consistent with definition of stochastic matrix ...
- That is why we divide by 2.

(19)



- Kopecky and Suen (2010, RES) show that the Rouwenhorst method is superior when the process is highly persistent
- Assume we have stochastic process

$$x_t = 0.975x_{t-1} + \epsilon$$
 with  $\mathcal{N}(0, 0.007^2)$  (21)

- Discretize the process using both methods, simulate using T = 10,000 and compute some moments

|       | Tauchen         |                  |        | Rouwenhorst     |                  |        |  |
|-------|-----------------|------------------|--------|-----------------|------------------|--------|--|
|       | $\hat{\mu}_{y}$ | $\hat{\sigma}_y$ | ρ      | $\hat{\mu}_{y}$ | $\hat{\sigma}_y$ | ρ      |  |
| n = 5 | 0.0009          | 0.0042           | 0.9969 | -0.001          | 0.0069           | 0.9753 |  |
| n = 9 | -0.0013         | 0.0076           | 0.9778 | 0.0003          | 0.0071           | 0.9763 |  |



- Tauchen method can be extended to VAR(1) processes
  - \* ARMA(p,q) can be written as VAR(1)
  - \* VAR(p) can also be written as VAR(1)
- What if the process is not stationary?
  - \* Typically the case of the income process in quantitative life-cycle models
  - \* Fella, Galliponi and Pan (2019, Rev. Econ. Dyn.)